

ESI

Ru/TiO₂-catalysed hydrogenation of xylose: the role of crystal structure of the support

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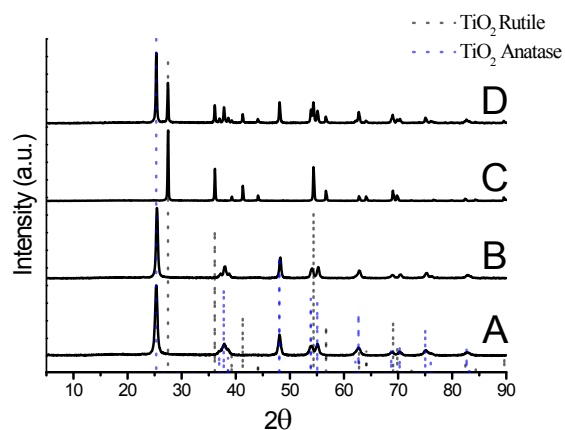


Figure S1. XRD patterns of the supports A–D.

Phase percent composition was determined by XRD using the following equation:

$$\%Rutile = \frac{1}{\left(\frac{A}{R} \times 0.884\right) + 1} \times 100$$

where A is the area of anatase major peak ($2\theta=25^\circ$), R the area of rutile major peak ($2\theta=28^\circ$) and 0.884 is a scattering coefficient¹.

Table S1: Product distribution at different reaction temperatures with each catalyst.

Catalysts	Temperature (°C)	% Conversion	%Yield					
			Xylitol	Arabitol	Adonitol	Furfural	Glycols	NI
A	120	70.7	10.4	2.3	0.0	1.4	17.4	39.3
	140	92.6	7.1	0.0	0.0	3.1	27.1	55.3
	160	95.6	4.6	3.0	0.0	5.9	43.0	39.1

B	120	84.9	65.2	0.9	2.3	0.0	6.7	9.8
	140	97.4	43.9	2.5	0.0	1.7	27.7	21.6
	160	96.9	34.1	0.9	0.0	0.0	34.9	26.9
C	120	100.0	93.6	5.4	0.9	0.0	0.0	0.0
	140	100.0	85.2	10.0	4.8	0.0	0.0	0.0
	160	100.0	70.8	16.0	6.8	0.0	0.0	6.4
D	120	100.0	95.5	2.5	2.0	0.0	0.0	0.0
	140	100.0	87.8	7.5	4.7	0.0	0.0	0.0
	160	100.0	72.2	18.8	9.0	0.0	0.0	0.0

NI: not identified.

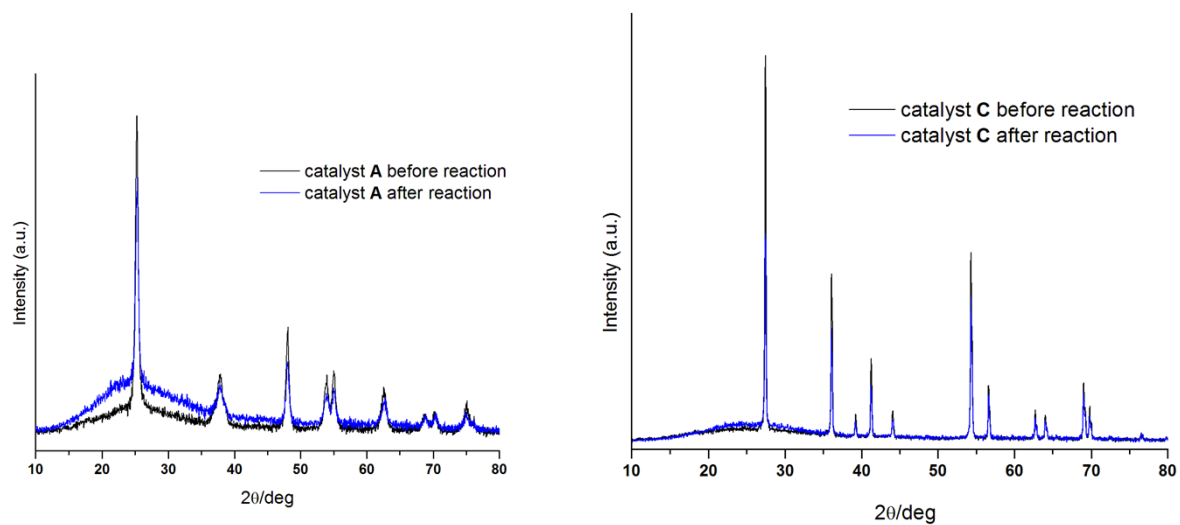
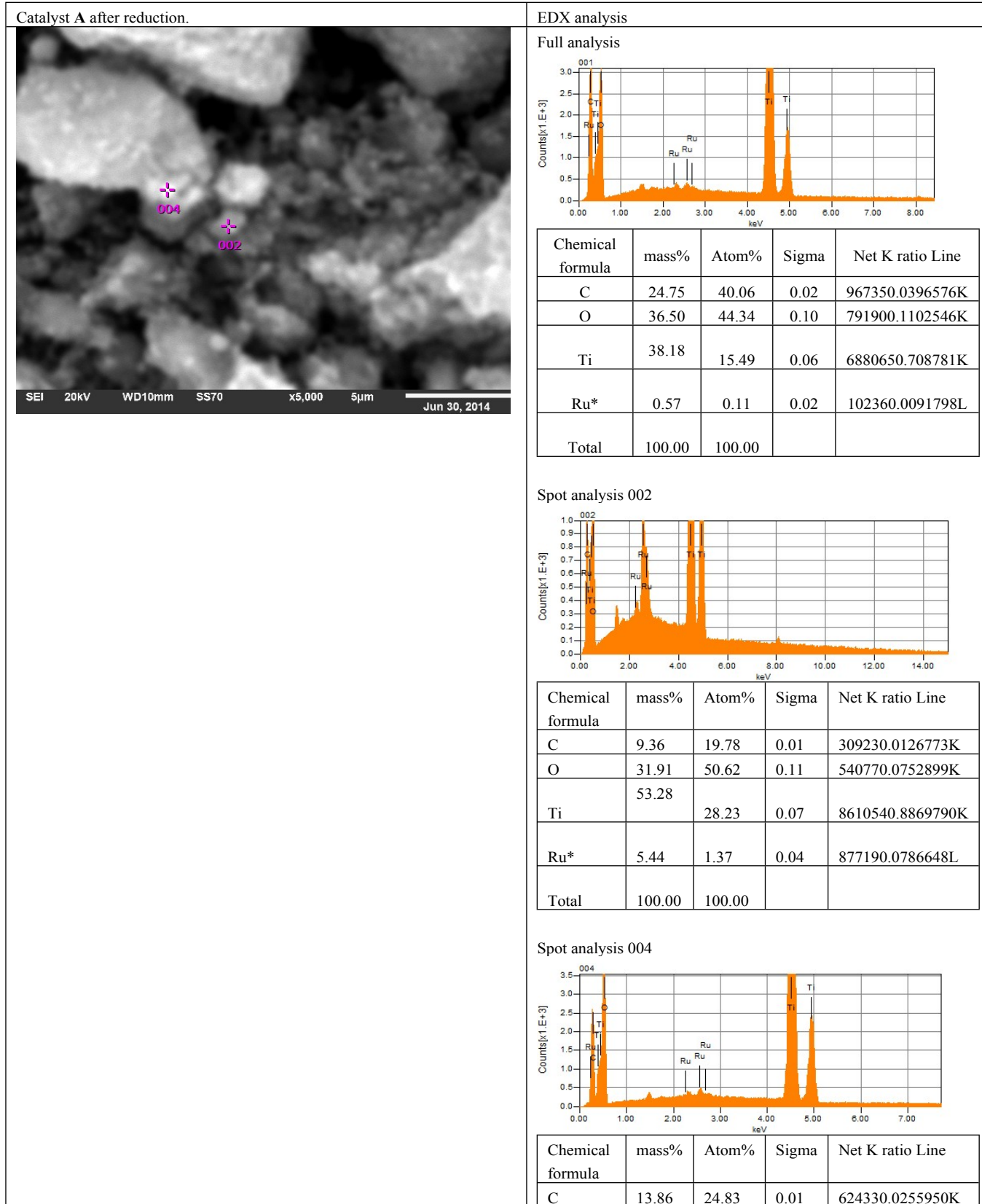
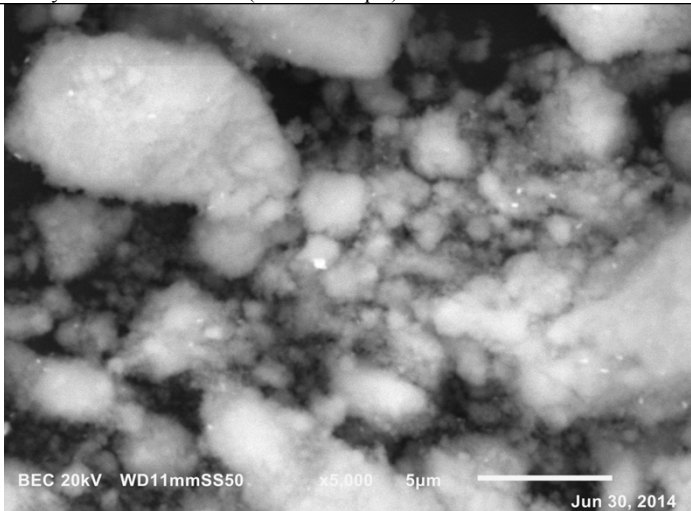
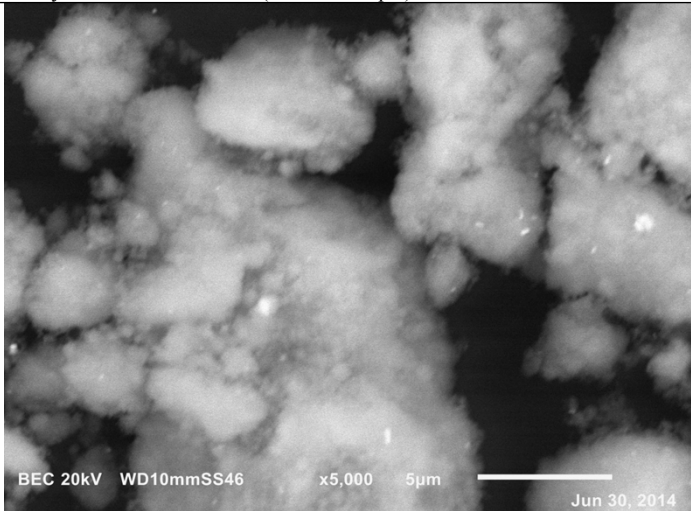
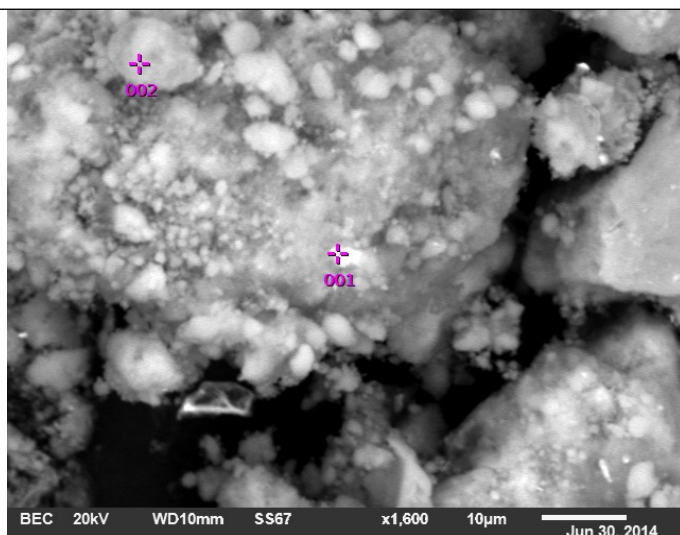


Fig. S2: XRD patterns of the catalysts A and C before and after reaction

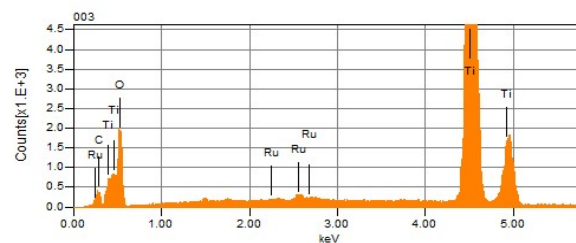
SEM and EDX analysis.



	<table><tr><td>O</td><td>40.83</td><td>54.92</td><td>0.09</td><td>1114280.1551384K</td></tr><tr><td>Ti</td><td>44.86</td><td>20.15</td><td>0.05</td><td>10068201.037133K</td></tr><tr><td>Ru*</td><td>0.46</td><td>0.10</td><td>0.02</td><td>102040.0091511L</td></tr><tr><td>Total</td><td>100.00</td><td>100.00</td><td></td><td></td></tr></table>				O	40.83	54.92	0.09	1114280.1551384K	Ti	44.86	20.15	0.05	10068201.037133K	Ru*	0.46	0.10	0.02	102040.0091511L	Total	100.00	100.00		
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Catalyst A after reduction. (BEC technique)																								
																								
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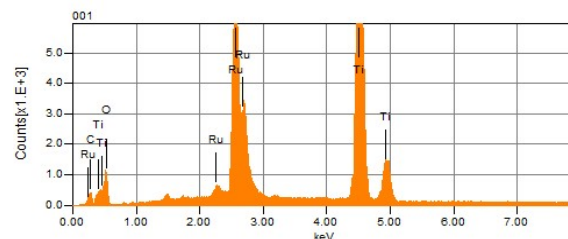


Full analysis.



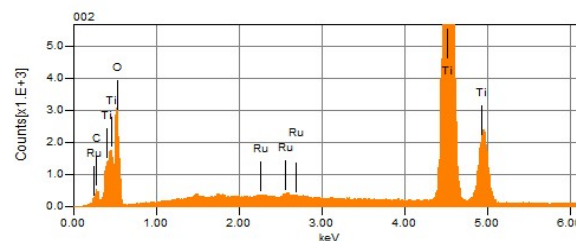
Chemical formula	mass%	Atom%	Sigma	Net K ratio Line
C	4.21	9.07	0.01	95370.0039100K
O	36.50	59.10	0.13	465840.0648572K
Ti	58.45	31.61	0.08	7013470.7224636K
Ru*	0.84	0.22	0.03	98850.0088642L
Total	100.00	100.00		

Spot analysis 001

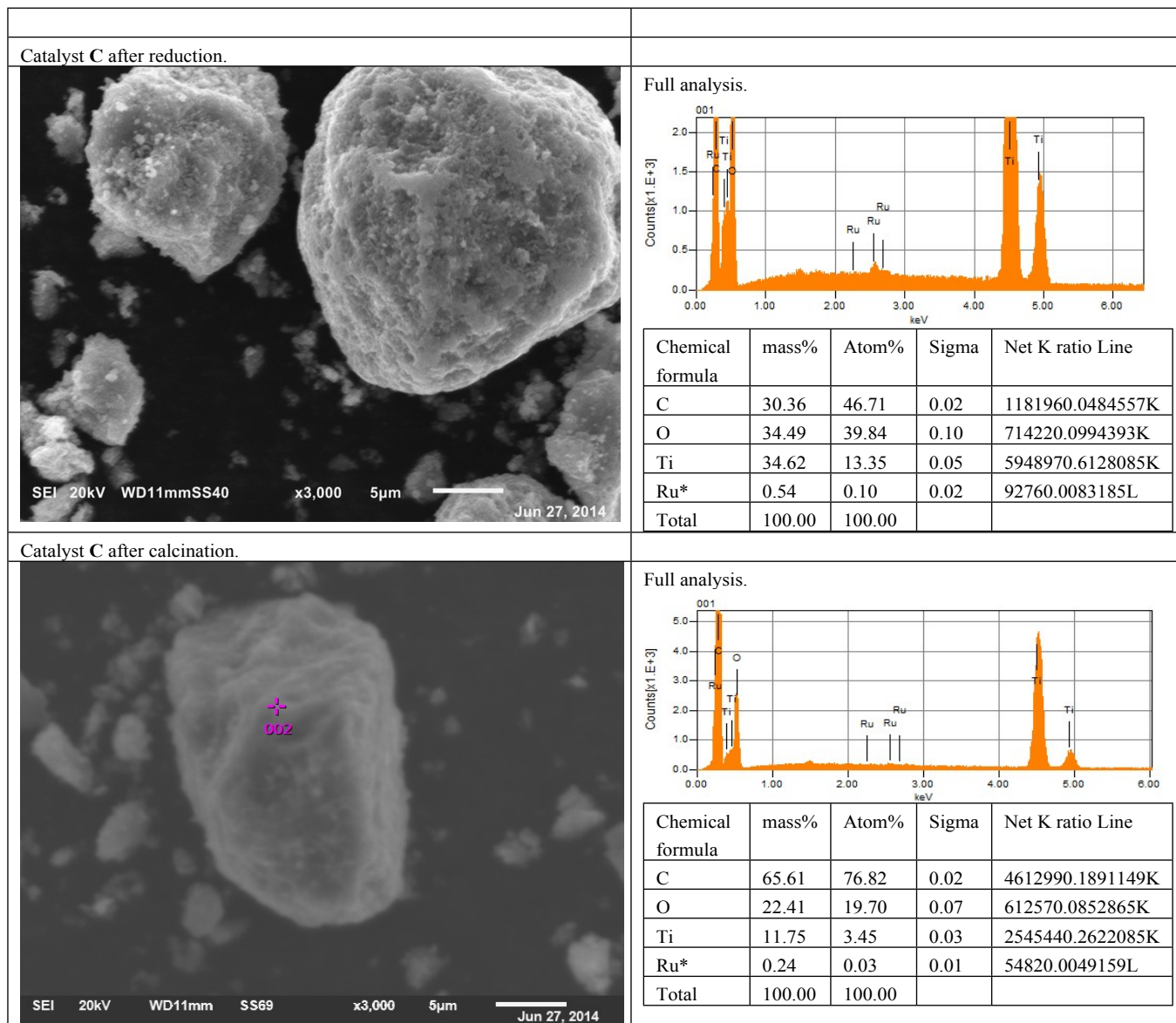


Chemical formula	mass%	Atom%	Sigma	Net K ratio Line
C	2.40	7.68	0.01	80600.0035619K
O	18.32	43.92	0.09	246520.0369980K
Ti	43.49	34.82	0.07	5295040.5879768K
Ru*	35.79	13.58	0.10	4707260.4550520L
Total	100.00	100.00		

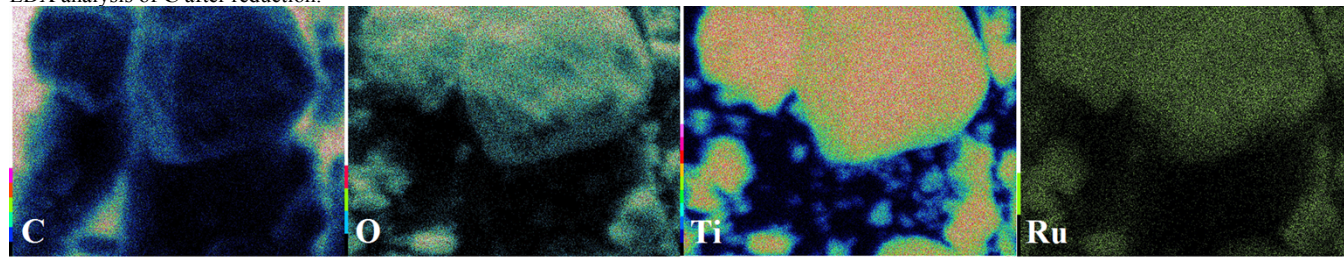
Spot analysis 002



Chemical formula	mass%	Atom%	Sigma	Net K ratio Line
C	3.43	7.31	0.01	105430.0043222K
O	38.75	61.94	0.12	696750.0970074K
Ti	57.39	30.64	0.07	9399680.9682691K
Ru*	0.43	0.11	0.03	6880 0.0061696 L
Total	100.00	100.00		

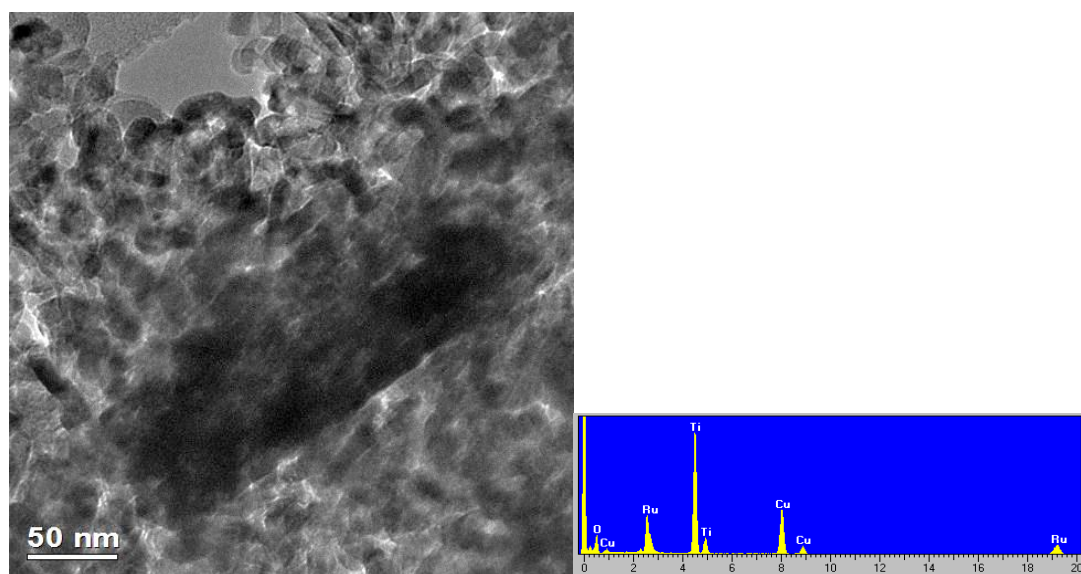
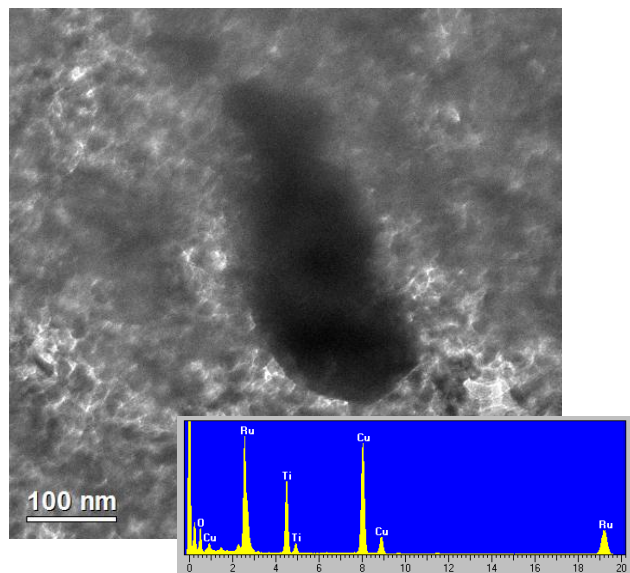


EDX analysis of C after reduction.



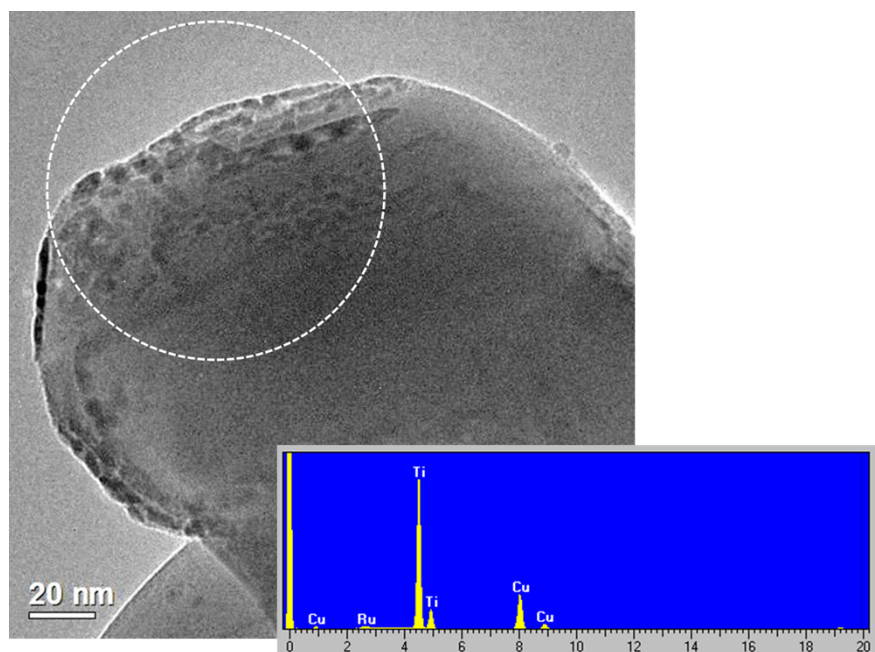
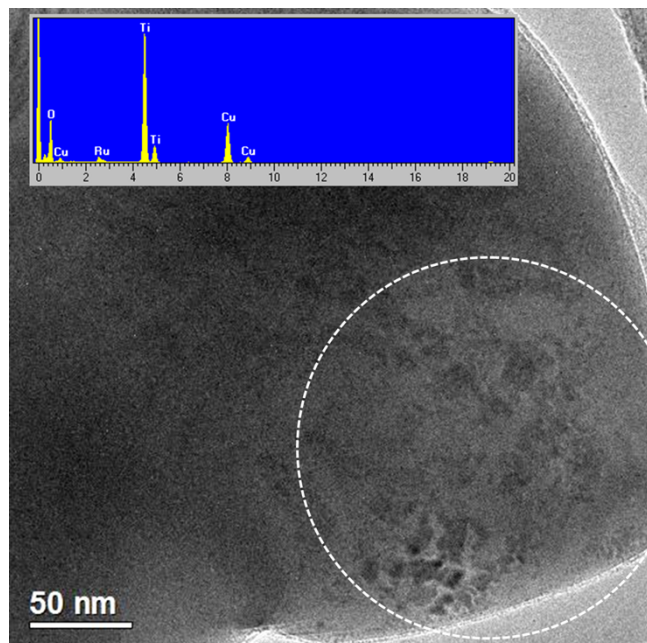
TEM analysis of spent samples

TEM images of two different regions of spent catalyst A. Corresponding EDX spectra were collected only from the dark contrasted Ru rich regions.



TEM images of two different regions of spent catalyst C.

EDX spectra from the selected areas (circled) show that Ru particles are small and highly dispersed after the reaction also.



1. X. Z. Fu, L. A. Clark, Q. Yang and M. A. Anderson, *Environmental Science & Technology*, 1996, 30, 647-653.